



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:51 PM GMT

PDB ID : 4JPV
Title : Crystal structure of broadly and potently neutralizing antibody 3bnc117 in complex with hiv-1 gp120
Authors : Zhou, T.; Moquin, S.; Kwong, P.D.
Deposited on : 2013-03-19
Resolution : 2.83 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

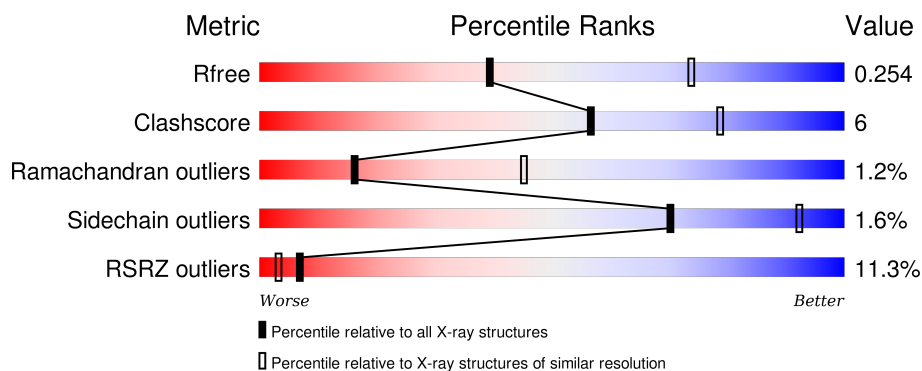
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.83 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	2676 (2.84-2.80)
Clashscore	102246	3124 (2.84-2.80)
Ramachandran outliers	100387	3072 (2.84-2.80)
Sidechain outliers	100360	3074 (2.84-2.80)
RSRZ outliers	91569	2690 (2.84-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	G	352	<div> <div>2%</div> <div>84%</div> <div>12%</div> <div>..</div> </div>
2	H	226	<div> <div>7%</div> <div>78%</div> <div>16%</div> <div>..</div> </div>
3	L	206	<div> <div>31%</div> <div>82%</div> <div>17%</div> <div>..</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	NAG	G	510	-	-	-	X
6	EDO	G	512	-	-	-	X
6	EDO	G	514	-	-	-	X
6	EDO	L	302	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6224 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called HIV-1 CLADE A STRAIN 93TH057 GP120 WITH LOOP d AND LOOPD V5 REPLACED FROM HIV STRAIN 3415V1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	G	342	Total	C	N	O	S	0	0	0
			2679	1683	464	510	22			

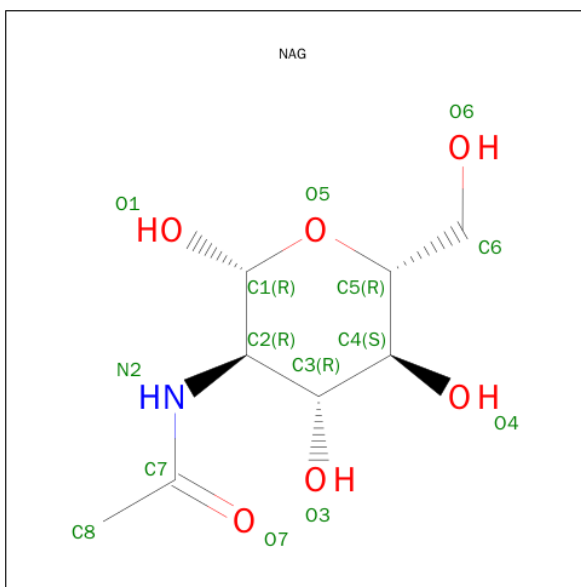
- Molecule 2 is a protein called HEAVY CHAIN OF ANTIBODY 3BNC117.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	216	Total	C	N	O	S	0	0	0
			1676	1065	289	317	5			

- Molecule 3 is a protein called LIGHT CHAIN OF ANTIBODY 3BNC117.

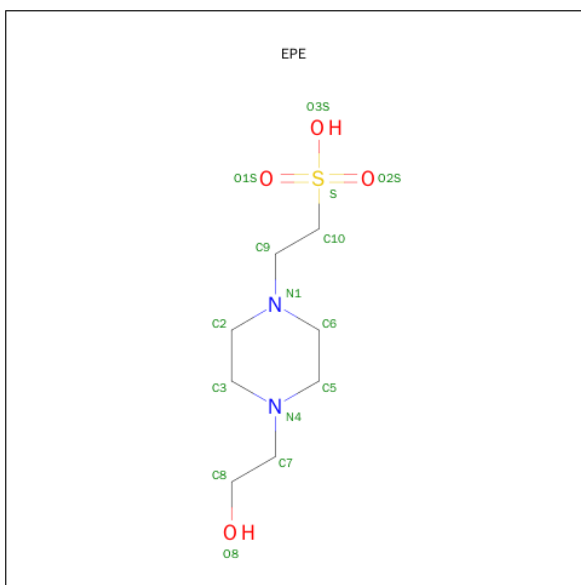
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	204	Total	C	N	O	S	0	0	0
			1603	1004	277	317	5			

- Molecule 4 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C₈H₁₅NO₆).



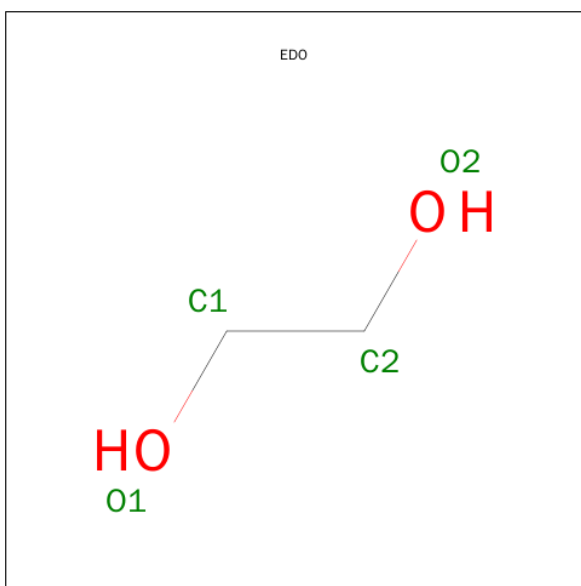
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	G	1	Total	C	N	O	0	0
			14	8	1	5		
4	G	1	Total	C	N	O	0	0
			14	8	1	5		
4	G	1	Total	C	N	O	0	0
			14	8	1	5		
4	G	1	Total	C	N	O	0	0
			14	8	1	5		
4	G	1	Total	C	N	O	0	0
			14	8	1	5		
4	G	1	Total	C	N	O	0	0
			14	8	1	5		
4	G	1	Total	C	N	O	0	0
			14	8	1	5		
4	L	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	G	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	G	1	Total	C	O	0	0
			4	2	2		
6	G	1	Total	C	O	0	0
			4	2	2		
6	G	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	L	1	Total	C	O	0	0
			4	2	2		

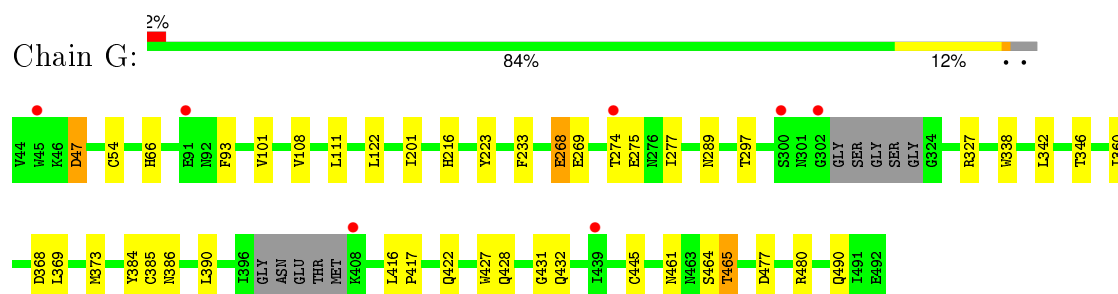
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	G	57	Total	O	0	0
			57	57		
7	H	11	Total	O	0	0
			11	11		
7	L	13	Total	O	0	0
			13	13		

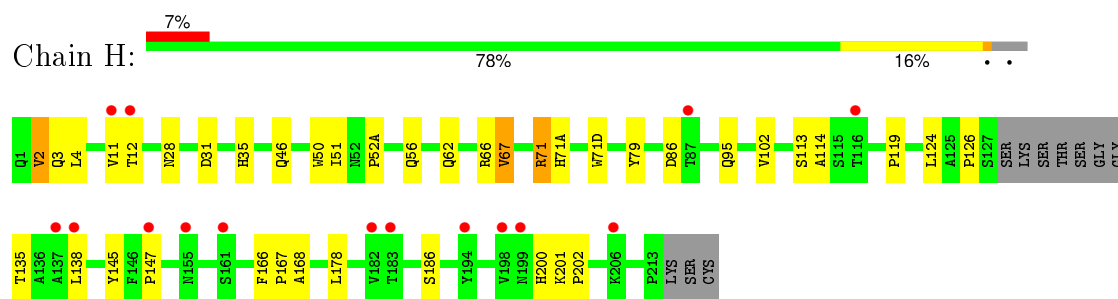
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

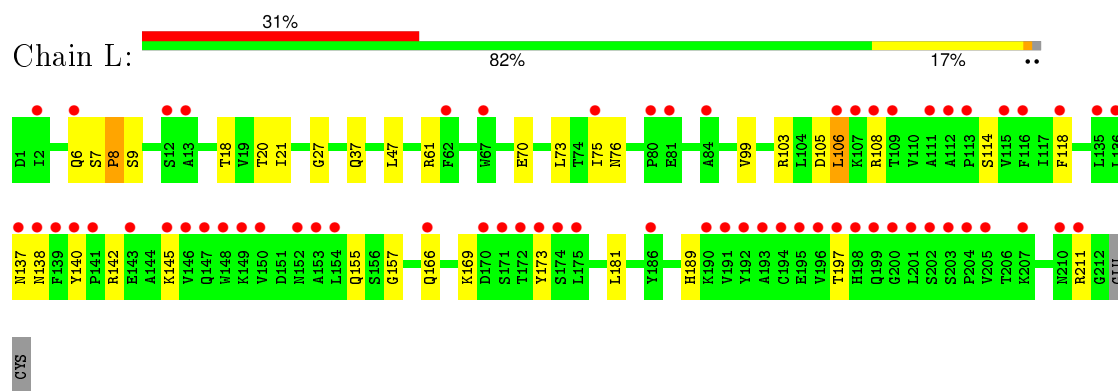
- Molecule 1: HIV-1 CLADE A STRAIN 93TH057 GP120 WITH LOOP d AND LOOPD V5 REPLACED FROM HIV STRAIN 3415V1



- Molecule 2: HEAVY CHAIN OF ANTIBODY 3BNC117



- Molecule 3: LIGHT CHAIN OF ANTIBODY 3BNC117



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.63Å 69.89Å 231.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.97 – 2.83 48.97 – 2.83	Depositor EDS
% Data completeness (in resolution range)	97.9 (48.97-2.83) 97.9 (48.97-2.83)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_998)	Depositor
R, R_{free}	0.199 , 0.259 0.192 , 0.254	Depositor DCC
R_{free} test set	1371 reflections (5.09%)	DCC
Wilson B-factor (Å ²)	68.2	Xtriage
Anisotropy	0.542	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 69.4	EDS
Estimated twinning fraction	0.038 for k,h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 26961 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6224	wwPDB-VP
Average B, all atoms (Å ²)	95.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.66% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: EPE, NAG, EDO

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	G	0.30	0/2735	0.50	0/3711
2	H	0.26	0/1726	0.45	0/2359
3	L	0.24	0/1637	0.42	0/2222
All	All	0.27	0/6098	0.47	0/8292

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	2679	0	2613	30	0
2	H	1676	0	1608	21	0
3	L	1603	0	1563	24	0
4	G	140	0	130	2	0
4	L	14	0	13	1	0
5	G	15	0	17	1	0
6	G	12	0	18	0	0
6	L	4	0	6	0	0
7	G	57	0	0	9	0
7	H	11	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	L	13	0	0	3	0
All	All	6224	0	5968	74	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (74) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:7:SER:O	3:L:9:SER:N	2.19	0.74
3:L:189:HIS:O	3:L:211:ARG:NH1	2.26	0.69
1:G:373:MET:HG2	1:G:384:TYR:HB3	1.77	0.67
3:L:7:SER:HB3	3:L:8:PRO:HD2	1.77	0.66
3:L:145:LYS:HB3	3:L:197:THR:HB	1.78	0.66
3:L:6:GLN:HB2	3:L:99:VAL:HG11	1.81	0.63
1:G:386:ASN:N	7:G:640:HOH:O	2.33	0.61
3:L:21:ILE:HD11	3:L:73:LEU:HD23	1.82	0.60
1:G:327:ARG:NH1	1:G:422:GLN:OE1	2.29	0.59
2:H:46:GLN:NE2	2:H:62:GLN:OE1	2.34	0.59
2:H:2:VAL:HG11	2:H:102:VAL:HG21	1.84	0.59
1:G:417:PRO:O	7:G:640:HOH:O	2.17	0.57
3:L:105:ASP:OD2	3:L:173:TYR:OH	2.15	0.57
2:H:119:PRO:HB3	2:H:145:TYR:HB3	1.87	0.56
1:G:368:ASP:OD1	2:H:71:ARG:NH2	2.38	0.55
2:H:126:PRO:HB3	2:H:138:LEU:HB3	1.88	0.54
2:H:11:VAL:HG21	2:H:147:PRO:HG3	1.89	0.54
3:L:105:ASP:OD1	3:L:106:LEU:N	2.41	0.54
1:G:66:HIS:CD2	1:G:111:LEU:HD21	2.43	0.53
2:H:66:ARG:NH2	2:H:86:ASP:OD2	2.41	0.53
1:G:122:LEU:HD21	1:G:431:GLY:HA3	1.91	0.52
2:H:168:ALA:HA	2:H:178:LEU:HB3	1.91	0.52
3:L:157:GLY:N	7:L:401:HOH:O	2.42	0.51
2:H:135:THR:N	2:H:186:SER:HG	2.09	0.50
1:G:428:GLN:N	1:G:428:GLN:OE1	2.40	0.50
1:G:360:ILE:HD12	1:G:465:THR:HG21	1.94	0.49
3:L:114:SER:HB2	3:L:137:ASN:HB3	1.95	0.49
1:G:268:GLU:HG3	1:G:289:ASN:HB2	1.95	0.49
1:G:432:GLN:OE1	7:G:605:HOH:O	2.20	0.48
1:G:47:ASP:N	1:G:47:ASP:OD1	2.46	0.48
3:L:6:GLN:HG3	3:L:7:SER:H	1.79	0.48
2:H:4:LEU:O	7:H:311:HOH:O	2.18	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:445:CYS:SG	7:G:648:HOH:O	2.61	0.47
2:H:28:ASN:ND2	2:H:31:ASP:OD2	2.47	0.47
3:L:70:GLU:OE2	4:L:301:NAG:O6	2.31	0.47
3:L:169:LYS:NZ	7:L:410:HOH:O	2.41	0.47
1:G:108:VAL:HG23	1:G:427:TRP:HH2	1.78	0.46
3:L:103:ARG:HH11	3:L:142:ARG:HH12	1.63	0.46
1:G:274:THR:OG1	1:G:277:ILE:HG12	2.15	0.46
3:L:37:GLN:HB2	3:L:47:LEU:HD11	1.97	0.45
1:G:223:TYR:CE2	1:G:490:GLN:HB2	2.52	0.45
4:G:504:NAG:H61	3:L:27:GLY:HA2	1.98	0.45
1:G:338:TRP:CZ2	1:G:390:LEU:HG	2.52	0.45
3:L:105:ASP:OD1	3:L:166:GLN:NE2	2.50	0.44
2:H:168:ALA:HB2	2:H:178:LEU:HD23	1.99	0.44
2:H:71(A):HIS:CD2	2:H:79:TYR:HE2	2.35	0.44
2:H:201:LYS:NZ	7:H:307:HOH:O	2.50	0.44
3:L:106:LEU:O	3:L:140:TYR:OH	2.26	0.43
2:H:52(A):PRO:O	2:H:71:ARG:HD2	2.18	0.43
1:G:54:CYS:HB3	7:G:626:HOH:O	2.19	0.43
2:H:200:HIS:CE1	2:H:202:PRO:HB2	2.54	0.42
2:H:35:HIS:CD2	2:H:95:GLN:HB2	2.54	0.42
1:G:101:VAL:HG21	1:G:480:ARG:HG2	2.01	0.42
1:G:416:LEU:HB3	7:G:640:HOH:O	2.18	0.42
2:H:124:LEU:HB3	3:L:118:PHE:CD1	2.53	0.42
3:L:20:THR:HA	3:L:73:LEU:O	2.20	0.42
5:G:511:EPE:H71	7:G:626:HOH:O	2.18	0.42
1:G:93:PHE:HB2	1:G:233:PHE:HZ	1.85	0.42
3:L:103:ARG:NH1	3:L:142:ARG:HH12	2.17	0.41
3:L:18:THR:HA	3:L:75:ILE:O	2.20	0.41
1:G:216:HIS:HE1	7:G:614:HOH:O	2.03	0.41
1:G:342:LEU:O	1:G:346:THR:HG23	2.21	0.41
1:G:477:ASP:OD1	1:G:480:ARG:NH1	2.53	0.41
2:H:51:ILE:HA	2:H:56:GLN:O	2.21	0.41
1:G:108:VAL:HG23	1:G:427:TRP:CH2	2.56	0.40
3:L:61:ARG:HB3	3:L:76:ASN:O	2.22	0.40
1:G:268:GLU:HA	1:G:269:GLU:HA	1.91	0.40
1:G:269:GLU:HB2	4:G:505:NAG:O6	2.21	0.40
1:G:373:MET:HG3	1:G:385:CYS:C	2.42	0.40
2:H:166:PHE:HA	2:H:167:PRO:HD3	1.98	0.40
3:L:108:ARG:HD2	7:L:405:HOH:O	2.21	0.40
1:G:461:ASN:O	1:G:464:SER:N	2.52	0.40
2:H:35:HIS:CE1	2:H:50:TRP:HB3	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:297:THR:HG23	7:G:617:HOH:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	G	336/352 (96%)	312 (93%)	22 (6%)	2 (1%)	30	63
2	H	212/226 (94%)	193 (91%)	14 (7%)	5 (2%)	7	24
3	L	202/206 (98%)	185 (92%)	15 (7%)	2 (1%)	19	51
All	All	750/784 (96%)	690 (92%)	51 (7%)	9 (1%)	16	45

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	275	GLU
2	H	113	SER
3	L	8	PRO
2	H	2	VAL
2	H	67	VAL
1	G	268	GLU
2	H	3	GLN
2	H	114	ALA
3	L	138	ASN

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	305/311 (98%)	301 (99%)	4 (1%)	76	94
2	H	185/193 (96%)	181 (98%)	4 (2%)	60	88
3	L	181/183 (99%)	178 (98%)	3 (2%)	68	91
All	All	671/687 (98%)	660 (98%)	11 (2%)	70	92

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	47	ASP
1	G	201	ILE
1	G	369	LEU
1	G	465	THR
2	H	12	THR
2	H	67	VAL
2	H	71	ARG
2	H	71(D)	TRP
3	L	106	LEU
3	L	155	GLN
3	L	181	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
3	L	155	GLN
3	L	166	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

16 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
4	NAG	G	501	1	14,14,15	0.49	0	15,19,21	0.79	0
4	NAG	G	502	1	14,14,15	0.51	0	15,19,21	1.10	1 (6%)
4	NAG	G	503	1	14,14,15	0.47	0	15,19,21	1.30	2 (13%)
4	NAG	G	504	1	14,14,15	0.60	0	15,19,21	0.81	0
4	NAG	G	505	1	14,14,15	0.56	0	15,19,21	1.01	1 (6%)
4	NAG	G	506	1	14,14,15	0.52	0	15,19,21	0.48	0
4	NAG	G	507	1	14,14,15	0.49	0	15,19,21	1.05	2 (13%)
4	NAG	G	508	1	14,14,15	0.62	0	15,19,21	0.75	0
4	NAG	G	509	1	14,14,15	0.54	0	15,19,21	0.81	0
4	NAG	G	510	1	14,14,15	0.48	0	15,19,21	0.81	0
5	EPE	G	511	-	14,15,15	0.36	0	18,20,20	1.80	3 (16%)
6	EDO	G	512	-	3,3,3	0.45	0	2,2,2	0.46	0
6	EDO	G	513	-	3,3,3	0.47	0	2,2,2	0.45	0
6	EDO	G	514	-	3,3,3	0.49	0	2,2,2	0.32	0
4	NAG	L	301	3	14,14,15	0.55	0	15,19,21	0.86	1 (6%)
6	EDO	L	302	-	3,3,3	0.52	0	2,2,2	0.30	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	G	501	1	-	0/6/23/26	0/1/1/1
4	NAG	G	502	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	G	503	1	-	0/6/23/26	0/1/1/1
4	NAG	G	504	1	-	0/6/23/26	0/1/1/1
4	NAG	G	505	1	-	0/6/23/26	0/1/1/1
4	NAG	G	506	1	-	0/6/23/26	0/1/1/1
4	NAG	G	507	1	-	0/6/23/26	0/1/1/1
4	NAG	G	508	1	-	0/6/23/26	0/1/1/1
4	NAG	G	509	1	-	0/6/23/26	0/1/1/1
4	NAG	G	510	1	-	0/6/23/26	0/1/1/1
5	EPE	G	511	-	-	0/9/19/19	0/1/1/1
6	EDO	G	512	-	-	0/1/1/1	0/0/0/0
6	EDO	G	513	-	-	0/1/1/1	0/0/0/0
6	EDO	G	514	-	-	0/1/1/1	0/0/0/0
4	NAG	L	301	3	-	0/6/23/26	0/1/1/1
6	EDO	L	302	-	-	0/1/1/1	0/0/0/0

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	503	NAG	C6-C5-C4	-2.46	106.94	113.02
4	G	507	NAG	O5-C5-C6	2.01	111.69	107.35
4	G	507	NAG	C1-O5-C5	2.08	114.89	112.25
4	G	505	NAG	C1-O5-C5	2.11	114.92	112.25
5	G	511	EPE	C7-N4-C3	2.24	117.02	111.27
4	L	301	NAG	C1-O5-C5	2.44	115.34	112.25
4	G	503	NAG	C1-O5-C5	3.40	116.57	112.25
4	G	502	NAG	C1-O5-C5	3.45	116.62	112.25
5	G	511	EPE	O2S-S-C10	3.85	110.19	106.91
5	G	511	EPE	C5-N4-C3	4.45	118.54	108.90

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	G	504	NAG	1	0
4	G	505	NAG	1	0
5	G	511	EPE	1	0
4	L	301	NAG	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	G	342/352 (97%)	0.32	7 (2%) 68 58	49, 70, 114, 164	0
2	H	216/226 (95%)	0.53	15 (6%) 20 11	50, 98, 157, 173	0
3	L	204/206 (99%)	1.40	64 (31%) 1 0	71, 128, 163, 208	0
All	All	762/784 (97%)	0.67	86 (11%) 7 3	49, 88, 158, 208	0

All (86) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	L	108	ARG	7.1
3	L	149	LYS	6.4
3	L	116	PHE	6.0
3	L	202	SER	5.9
3	L	200	GLY	5.8
3	L	112	ALA	5.2
3	L	201	LEU	5.0
3	L	111	ALA	4.9
3	L	195	GLU	4.8
3	L	199	GLN	4.6
3	L	172	THR	4.2
3	L	115	VAL	4.1
3	L	113	PRO	4.1
3	L	204	PRO	4.1
3	L	191	VAL	4.1
3	L	203	SER	4.1
3	L	190	LYS	4.1
2	H	206	LYS	4.0
3	L	150	VAL	4.0
3	L	205	VAL	3.9
3	L	153	ALA	3.9
2	H	137	ALA	3.7
3	L	194	CYS	3.7

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Mol	Chain	Res	Type	RSRZ
3	L	106	LEU	3.7
2	H	12	THR	3.7
2	H	182	VAL	3.6
3	L	140	TYR	3.6
3	L	109	THR	3.6
3	L	107	LYS	3.6
3	L	135	LEU	3.4
3	L	139	PHE	3.3
3	L	118	PHE	3.2
3	L	210	ASN	3.2
3	L	198	HIS	3.2
2	H	138	LEU	3.2
3	L	12	SER	3.1
3	L	148	TRP	3.1
3	L	6	GLN	3.0
3	L	154	LEU	3.0
3	L	143	GLU	3.0
3	L	137	ASN	3.0
3	L	196	VAL	2.9
3	L	67	TRP	2.9
3	L	62	PHE	2.9
3	L	13	ALA	2.8
3	L	75	ILE	2.8
2	H	199	ASN	2.8
2	H	87	THR	2.7
2	H	198	VAL	2.7
2	H	161	SER	2.7
3	L	147	GLN	2.7
3	L	197	THR	2.6
3	L	175	LEU	2.5
3	L	145	LYS	2.5
3	L	192	TYR	2.5
1	G	408	LYS	2.5
1	G	91	GLU	2.4
3	L	170	ASP	2.4
3	L	166	GLN	2.4
3	L	2	ILE	2.4
3	L	152	ASN	2.3
2	H	147	PRO	2.3
1	G	300	SER	2.3
3	L	80	PRO	2.3
2	H	155	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
2	H	116	THR	2.3
3	L	138	ASN	2.3
3	L	173	TYR	2.2
3	L	211	ARG	2.2
3	L	207	LYS	2.2
1	G	274	THR	2.2
3	L	171	SER	2.1
2	H	183	THR	2.1
3	L	193	ALA	2.1
2	H	194	TYR	2.1
3	L	81	GLU	2.1
2	H	11	VAL	2.1
3	L	141	PRO	2.1
3	L	146	VAL	2.1
3	L	84	ALA	2.1
1	G	45	TRP	2.0
3	L	136	LEU	2.0
3	L	174	SER	2.0
1	G	302	GLY	2.0
1	G	439	ILE	2.0
3	L	186	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
6	EDO	G	512	4/4	0.94	0.48	9.08	76,79,79,80	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
6	EDO	L	302	4/4	0.71	0.37	7.63	88,88,91,92	0
4	NAG	G	510	14/15	0.76	0.43	4.92	82,110,120,126	0
6	EDO	G	514	4/4	0.86	0.28	3.65	77,79,84,86	0
4	NAG	G	502	14/15	0.93	0.23	0.74	66,87,102,104	0
4	NAG	G	507	14/15	0.94	0.21	-0.12	54,63,74,80	0
4	NAG	G	506	14/15	0.96	0.20	-0.73	46,69,84,94	0
4	NAG	G	503	14/15	0.95	0.18	-0.80	37,49,59,86	0
4	NAG	L	301	14/15	0.85	0.22	-0.80	104,122,138,139	0
4	NAG	G	501	14/15	0.94	0.16	-0.96	52,76,92,100	0
5	EPE	G	511	15/15	0.98	0.18	-1.28	33,45,70,79	0
4	NAG	G	505	14/15	0.97	0.17	-1.47	48,65,86,100	0
4	NAG	G	504	14/15	0.94	0.15	-5.56	74,90,99,101	0
6	EDO	G	513	4/4	0.78	0.19	-	97,104,105,105	0
4	NAG	G	508	14/15	0.95	0.17	-	66,76,86,90	0
4	NAG	G	509	14/15	0.88	0.22	-	119,132,139,141	0

6.5 Other polymers [i](#)

There are no such residues in this entry.